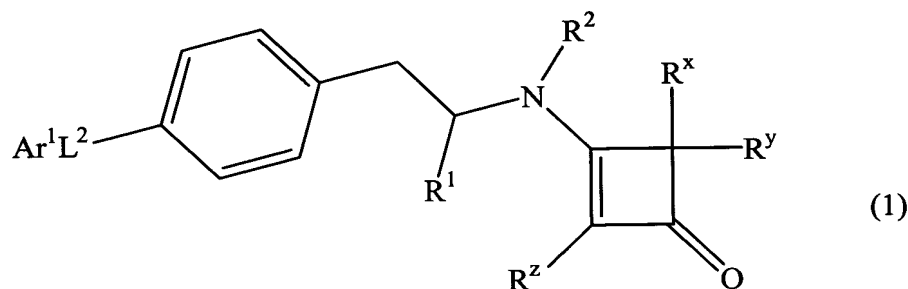


This listing of claims will replace all prior versions, and listings, of claims in the application.

Listing of Claims:

1. (currently amended) A process for the preparation of a compound of formula (1):



wherein:

Ar¹ is an optionally substituted aromatic or heteroaromatic group;

L² is a linker group selected from -N(R⁴)- [~~where R⁴ is a hydrogen atom or an optionally substituted straight or branched C₁₋₆alkyl group~~], -CON(R⁴)- and -S(O)₂N(R⁴)-;

R⁴ is a hydrogen atom or an optionally substituted straight or branched C₁₋₆alkyl group;

R¹ is a carboxylic acid (-CO₂H) or a derivative or biostere thereof an acyclic or cyclic carboxylic acid ester, an amide, tetrazole, phosphonic acid, phosphinic acid, sulphonic acid, sulphinic acid, boronic acid, or an acylsulphonamide group;

R² is a hydrogen atom or a C₁₋₆alkyl group;

R^x, R^y and R^z, which may be the same or different, are each an atom or group -L¹(Alk¹)_n(R³)_v in which, or R^z is -L¹(Alk¹)_n(R³)_v and R^x and R^y are joined together to form an optionally substituted spiro linked cycloaliphatic or heterocycloaliphatic group;

L¹ is a covalent bond or a linker atom or group an -O-, -S-, or -Se- atom or an -C(O)-, -C(O)O-, -OC(O)-, -C(S)-, -S(O)-, -S(O)₂-, -N(R⁸)-, -CON(R⁸)-, -OC(O)N(R⁸)-, -CSN(R⁸), -N(R⁸)CO-, -N(R⁸)C(O)O-, -N(R⁸)CS-, -S(O)₂N(R⁸)-, -N(R⁸)S(O)₂-, -N(R⁸)O-, -ON(R⁸)-, -N(R⁸)N(R⁸)-, -N(R⁸)CON(R⁸)-, -N(R⁸)CSN(R⁸)-, or -N(R⁸)SO₂N(R⁸)- group;

R⁸ is a hydrogen atom or an optionally substituted straight or branched C₁₋₆alkyl group;

Alk¹ is an optionally substituted aliphatic chain or an optionally substituted heteroaliphatic chain containing one to four -O- or -S- atoms or -C(O)-, -C(O)O-, -OC(O)-, -C(S)-, -S(O)-, -S(O)₂-, -N(R⁸)-, -CON(R⁸)-, -OC(O)N(R⁸)-, -CSN(R⁸), -N(R⁸)CO-, -N(R⁸)C(O)O-, -N(R⁸)CS-, -S(O)₂N(R⁸)-, -N(R⁸)S(O)₂-, -N(R⁸)O-, -ON(R⁸)-, -N(R⁸)N(R⁸)-, -N(R⁸)CON(R⁸)-, -N(R⁸)CSN(R⁸)-, or -N(R⁸)SO₂N(R⁸)- groups that interrupt or are at the terminus of the aliphatic chain;

R³ is a hydrogen or halogen atom or group selected from ~~-OR^{3a} [where R^{3a} is a hydrogen atom or an optionally substituted straight or branched C₁₋₆alkyl group or C₃₋₈cycloalkyl group]~~, -SR^{3a}, -CN and an optionally substituted cycloaliphatic, heterocycloaliphatic, polycycloaliphatic, heteropolycycloaliphatic, aromatic or heteroaromatic group;

R^{3a} is a hydrogen atom or an optionally substituted straight or branched C₁₋₆alkyl group or C₃₋₈cycloalkyl group;

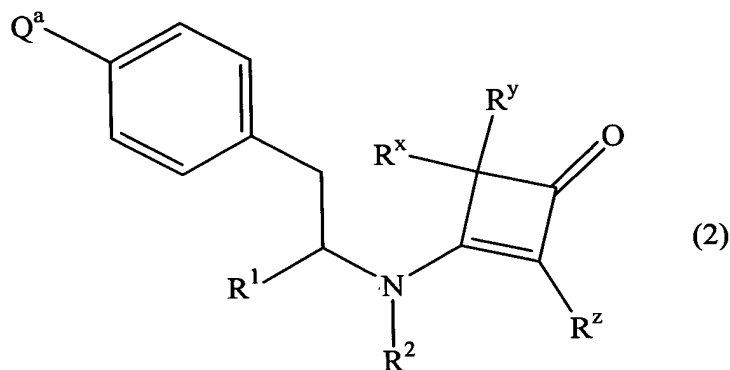
n is zero or the integer 1; and

v is the integer 1, 2 or 3;

provided that when n is zero and L¹ is a covalent bond, v is the integer 1;

~~or R^Z is an atom or group as previously defined and R^X and R^Y are joined together to form an optionally substituted spiro-linked cycloaliphatic or heterocycloaliphatic group;~~
and the salts, solvates, hydrates and N-oxides thereof;

which comprises reacting a compound of formula (2):



wherein:

Q^a is a group $-N(R^4)H$;

and the salts, solvates, hydrates and N-oxides thereof;

with a compound Ar^1W wherein

W is a group selected from X^1 (~~wherein X^1 is a leaving atom or group~~), $-COX^2$

(~~wherein X^2 is a halogen atom or a -OH group~~) and

$-SO_2X^3$ (~~in which X^3 is a halogen atom~~);

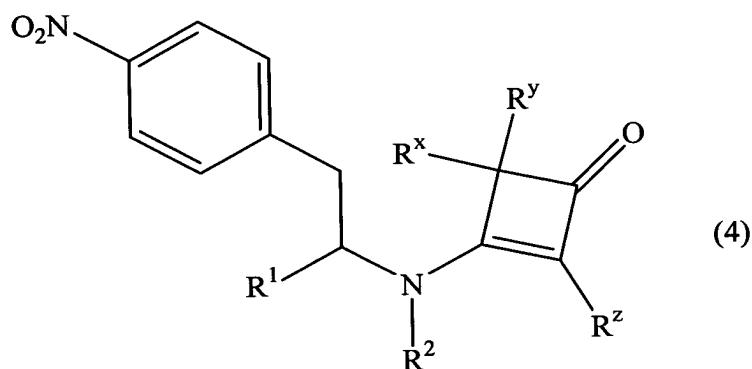
X^1 is a leaving atom or group;

X^2 is a halogen atom or a -OH group; and

X^3 is a halogen atom.

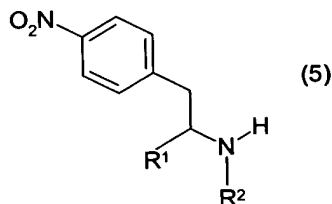
2. (original) A process according to Claim 1 wherein the reaction is carried out in a solvent in the presence of an acid when W is the group X^1 .
3. (previously presented) A process according to Claim 2 wherein the solvent is selected from an alcohol, ether, acetic acid, water, acetonitrile, substituted amide and ester.
4. (original) A process according to Claim 2 wherein the reaction is carried out in an alcohol in the presence of an acid catalyst.

5. (original) A process according to Claim 1 wherein the reaction is carried out in the presence of a base, an organic amine or a cyclic amine and an organic solvent when W is the group COX^2 and X^2 is a halogen atom.
6. (previously presented) A process according to Claim 5 wherein the organic solvent is selected from a halogenated hydrocarbon, a dipolar aprotic solvent, an ether and an ester.
7. (original) A process according to Claim 1 wherein the reaction is carried out in the presence of a condensing agent and a halogenated hydrocarbon, dipolar aprotic or an ether solvent when W is the group CO_2H .
8. (original) A process according to Claim 1 wherein the reaction is carried out in the presence of a base, an organic amine or a cyclic amine and a halogenated hydrocarbon, dipolar aprotic or an ether solvent when W is the group SO_2X^3 .
9. (previously presented) A process according to claim 1 wherein the compound of formula (2) is prepared by reduction of a compound of formula (4):

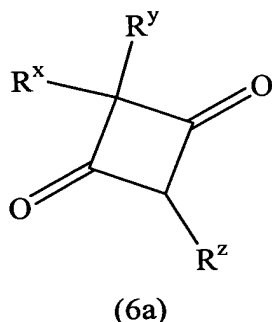


10. (original) A process according to Claim 9 wherein the reduction is carried out by catalytic hydrogenation or by chemical reduction.

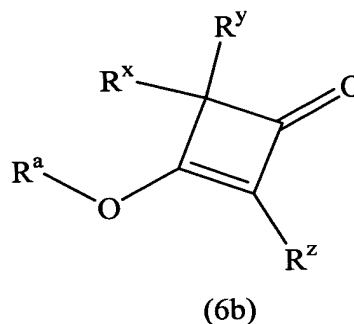
11. (previously presented) A process according to Claim 1 wherein R^4 is a hydrogen atom.
12. (original) A process according to Claim 9 wherein the compound of formula (4) is prepared by reaction of a compound of formula (5):



with a compound of formula (6a) or (6b):



or



wherein R^a represents a C₁-6alkyl group or a silyl group.

13. (original) A process according to Claim 12 wherein the reaction is carried out in the presence of an organic solvent.
14. (previously presented) A process according to Claim 13 wherein the solvent is selected from an aromatic hydrocarbon, a halogenated hydrocarbon and an ester.
15. (currently amended) A process according to Claim 1 wherein R^1 is the group $-\text{CO}_2\text{Alk}^7$; and

Alk^7 is a straight or branched optionally substituted C₁-8alkyl group, an optionally substituted C₂-8alkenyl group, an optionally substituted C₂-8alkynyl group, an optionally substituted C₃-8cycloalkyl group, an optionally substituted C₃-8heterocycloalkyl group, an

optionally substituted C₃-8cycloalkylC₁-8alkyl group, an optionally substituted C₃-8heterocycloalkylC₁-8alkyl group, an optionally substituted C₁-6alkyloxyC₁-6alkyl group, an optionally substituted hydroxyC₁-6alkyl group, an optionally substituted C₁-6alkylthioC₁-6alkyl group, an optionally substituted C₁-6alkylsulfinylC₁-6alkyl group, an optionally substituted C₁-6alkylsulfonylC₁-6alkyl group, an optionally substituted C₃-8cycloalkyloxyC₁-6alkyl group, an optionally substituted C₃-8cycloalkylthioC₁-6alkyl group, an optionally substituted C₃-8cycloalkylsulfinylC₁-6alkyl group, an optionally substituted C₃-8cycloalkylsulfonylC₁-6alkyl group, an optionally substituted C₁-6alkyloxycarbonylC₁-6alkyl group, an optionally substituted C₁-6alkyloxycarbonylC₁-6alkenyl group, an optionally substituted C₁-6alkyloxycarbonyloxyC₁-6alkyl group, an optionally substituted C₁-6alkyloxycarbonyloxyC₁-6alkenyl group, an optionally substituted C₃-8cycloalkyloxycarbonyloxyC₁-6alkyl group, an optionally substituted N-di-C₁-8alkylaminoC₁-8alkyl group, an optionally substituted N-C₆-12aryl-N-C₁-6alkylaminoC₁-6alkyl group, an optionally substituted N-di-C₁-8alkyl-carbamoylC₁-8alkyl group, an optionally substituted C₆-12arylC₁-6alkyl group, an optionally substituted heteroC₆-10arylC₁-6alkyl group, an optionally substituted C₆-12aryl group, an optionally substituted C₆-12aryloxyC₁-8alkyl group, an optionally substituted C₆-12arylthioC₁-8alkyl group, an optionally substituted C₆-12arylsulfinylC₁-8alkyl group, an optionally substituted C₆-12arylsulfonylC₁-8alkyl group, an optionally substituted C₁-8alkanoyloxyC₁-8alkyl group, an optionally substituted C₄-8imidoC₁-8alkyl group, an optionally substituted C₆-12aroyloxyC₁-8alkyl group, or a triglyceride.

16. (canceled)

17. (currently amended) A process according to ~~Claim 16~~ Claim 1 which comprises hydrolysing a compound of formula (1) in which R¹ is -CO₂Alk⁷ and Alk⁷ is a straight or branched optionally substituted C₁-8alkyl group, an optionally substituted C₂-8alkenyl group, an optionally substituted C₂-8alkynyl group, an optionally substituted C₃-8cycloalkyl

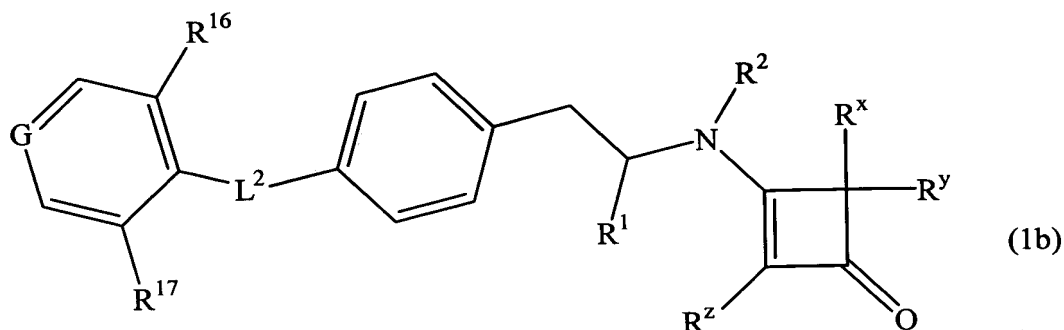
group, an optionally substituted C₃₋₈heterocycloalkyl group, an optionally substituted C₃₋₈cycloalkylC₁₋₈alkyl group, an optionally substituted C₃₋₈heterocycloalkylC₁₋₈alkyl group, an optionally substituted C₁₋₆alkyloxyC₁₋₆alkyl group, an optionally substituted hydroxyC₁₋₆alkyl group, an optionally substituted C₁₋₆alkylthioC₁₋₆alkyl group, an optionally substituted C₁₋₆alkylsulfinylC₁₋₆alkyl group, an optionally substituted C₁₋₆alkylsulfonylC₁₋₆alkyl group, an optionally substituted C₃₋₈cycloalkyloxyC₁₋₆alkyl group, an optionally substituted C₃₋₈cycloalkylthioC₁₋₆alkyl group, an optionally substituted C₃₋₈cycloalkylsulfinylC₁₋₆alkyl group, an optionally substituted C₃₋₈cycloalkylsulfonylC₁₋₆alkyl group, an optionally substituted C₁₋₆alkyloxycarbonylC₁₋₆alkyl group, an optionally substituted C₁₋₆alkyloxycarbonylC₁₋₆alkenyl group, an optionally substituted C₁₋₆alkyloxycarbonyloxyC₁₋₆alkyl group, an optionally substituted C₁₋₆alkyloxycarbonyloxyC₁₋₆alkenyl group, an optionally substituted C₃₋₈cycloalkyloxycarbonyloxyC₁₋₆alkyl group, an optionally substituted N-di-C₁₋₈alkylaminoC₁₋₈alkyl group, an optionally substituted N-C₆₋₁₂aryl-N-C₁₋₆alkylaminoC₁₋₆alkyl group, an optionally substituted N-di-C₁₋₈alkyl-carbamoylC₁₋₈alkyl group, an optionally substituted C₆₋₁₂arylC₁₋₆alkyl group, an optionally substituted heteroC₆₋₁₀arylC₁₋₆alkyl group, an optionally substituted C₆₋₁₂aryl group, an optionally substituted C₆₋₁₂aryloxyC₁₋₈alkyl group, an optionally substituted C₆₋₁₂arylthioC₁₋₈alkyl group, an optionally substituted C₆₋₁₂arylsulfinylC₁₋₈alkyl group, an optionally substituted C₆₋₁₂arylsulfonylC₁₋₈alkyl group, an optionally substituted C₁₋₈alkanoyloxyC₁₋₈alkyl group, an optionally substituted C₄₋₈imidoC₁₋₈alkyl group, an optionally substituted C₆₋₁₂aroyloxyC₁₋₈alkyl group, or a triglyceride,

to produce a compound of formula (1) in which R¹ is -CO₂H.

18. (currently amended) A process according to ~~Claim 16~~ Claim 1 which comprises esterifying a compound of formula (1) in which R¹ is -CO₂H to produce a compound of formula (1) in which R¹ is -CO₂Alk⁷ and Alk⁷ is a straight or branched optionally substituted C₁₋₈alkyl group, an optionally substituted C₂₋₈alkenyl group, an optionally substituted C₂₋

galkynyl group, an optionally substituted C₃-gcycloalkyl group, an optionally substituted C₃-gheterocycloalkyl group, an optionally substituted C₃-gcycloalkylC₁-galkyl group, an optionally substituted C₃-gheterocycloalkylC₁-galkyl group, an optionally substituted C₁-6alkyloxyC₁-6alkyl group, an optionally substituted hydroxyC₁-6alkyl group, an optionally substituted C₁-6alkylthioC₁-6alkyl group, an optionally substituted C₁-6alkylsulfinylC₁-6alkyl group, an optionally substituted C₁-6alkylsulfonylC₁-6alkyl group, an optionally substituted C₃-gcycloalkyloxyC₁-6alkyl group, an optionally substituted C₃-gcycloalkylthioC₁-6alkyl group, an optionally substituted C₃-gcycloalkylsulfinylC₁-6alkyl group, an optionally substituted C₃-gcycloalkylsulfonylC₁-6alkyl group, an optionally substituted C₁-6alkyloxycarbonylC₁-6alkyl group, an optionally substituted C₁-6alkyloxycarbonylC₁-6alkenyl group, an optionally substituted C₁-6alkyloxycarbonyloxyC₁-6alkyl group, an optionally substituted C₁-6alkyloxycarbonyloxyC₁-6alkenyl group, an optionally substituted C₃-gcycloalkyloxycarbonyloxyC₁-6alkyl group, an optionally substituted N-di-C₁-galkylaminoC₁-galkyl group, an optionally substituted N-C₆-12aryl-N-C₁-6alkylaminoC₁-6alkyl group, an optionally substituted N-di-C₁-galkyl-carbamoylC₁-galkyl group, an optionally substituted C₆-12arylC₁-6alkyl group, an optionally substituted heteroC₆-10arylC₁-6alkyl group, an optionally substituted C₆-12aryl group, an optionally substituted C₆-12aryloxyC₁-galkyl group, an optionally substituted C₆-12arylthioC₁-galkyl group, an optionally substituted C₆-12arylsulfinylC₁-galkyl group, an optionally substituted C₆-12arylsulfonylC₁-galkyl group, an optionally substituted C₁-galkanoyloxyC₁-galkyl group, an optionally substituted C₄-gimidoC₁-galkyl group, an optionally substituted C₆-12aroyloxyC₁-galkyl group, or a triglyceride.

19. (currently amended) A process according to Claim 1 for the preparation of compounds of formula (1b):



wherein

$-G=$ is $-CR^{18}=$, $-N=$ or $-N(O)=$;

R^{16} , R^{17} and R^{18} , which may be the same or different, are each a hydrogen atom or an atom or group $-L^3(Alk^2)_tL^4(R^5)_u$;

L^3 and L^4 are, independently, a covalent bond, an $-O-$ or $-S-$ atom, or a $-C(O)-$, $-C(O)O-$, $-OC(O)-$, $-C(S)-$, $-S(O)-$, $-S(O)_2-$, $-N(R^8)-$, $-CON(R^8)-$, $-OC(O)N(R^8)-$, $-CSN(R^8)-$, $-N(R^8)CO-$, $-N(R^8)C(O)O-$, $-N(R^8)CS-$, $-S(O)_2N(R^8)-$, $-N(R^8)S(O)_2-$, $-N(R^8)O-$, $-ON(R^8)-$, $-N(R^8)N(R^8)-$, $-N(R^8)CON(R^8)-$, $-N(R^8)CSN(R^8)-$, or $-N(R^8)SO_2N(R^8)-$ group;

R^8 is a hydrogen atom or an optionally substituted straight or branched C_{1-6} alkyl group;

t is zero or the integer 1;

u is an integer 1, 2 or 3;

Alk^2 is an optionally substituted aliphatic chain or an optionally substituted heteroaliphatic chain containing one to four $-O-$ or $-S-$ atoms or $-C(O)-$, $-C(O)O-$, $-OC(O)-$, $-C(S)-$, $-S(O)-$, $-S(O)_2-$, $-N(R^8)-$, $-CON(R^8)-$, $-OC(O)N(R^8)-$, $-CSN(R^8)-$, $-N(R^8)CO-$, $-N(R^8)C(O)O-$, $-N(R^8)CS-$, $-S(O)_2N(R^8)-$, $-N(R^8)S(O)_2-$, $-N(R^8)O-$, $-ON(R^8)-$, $-N(R^8)N(R^8)-$, $-N(R^8)CON(R^8)-$, $-N(R^8)CSN(R^8)-$, or $-N(R^8)SO_2N(R^8)-$ groups that interrupt or are at the terminus of the aliphatic chain;

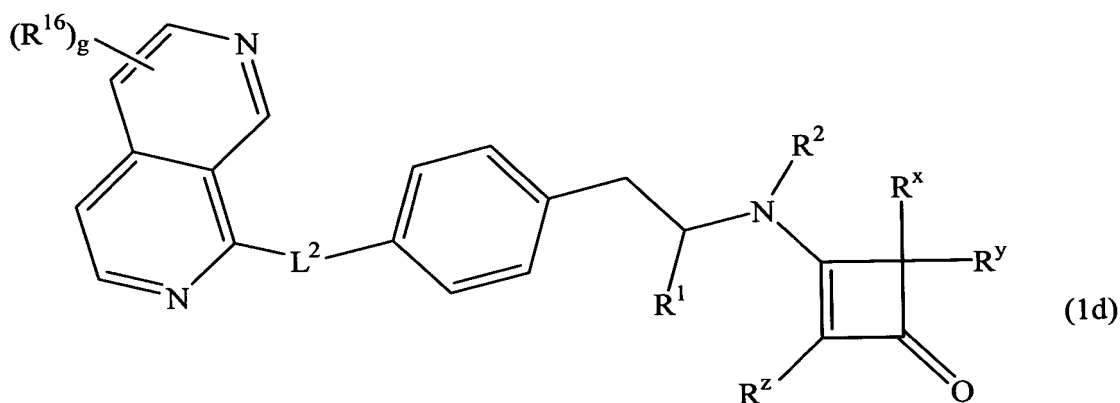
R^5 is a hydrogen or halogen atom or an optionally substituted C_{1-6} alkyl, optionally substituted C_{3-8} cycloalkyl, $-OR^6$, $-SR^6$, $-NR^6R^7$, $-NO_2$, $-CN$, $-CO_2R^6$, $-SO_3H$, $-SOR^6$, $-SO_2R^6$, $-SO_3R^6$, $-OCO_2R^6$, $-CONR^6R^7$, $-OCONR^6R^7$, $-CSNR^6R^7$, $-COR^6$, $-OCOR^6$,

-N(R⁶)COR⁷, -N(R⁶)CSR⁷, -SO₂N(R⁶)(R⁷), -N(R⁶)SO₂R⁷, N(R⁶)CON(R⁷)(R¹⁹),
 -N(R⁶)CSN(R⁷)(R¹⁹), or -N(R⁶)SO₂N(R⁷)(R¹⁹) group; and

R⁶, R⁷, and R¹⁹ are, independently, a hydrogen atom or an optionally substituted C₁-
 6alkyl or C₃-8cycloalkyl group;

provided that when t is zero and each of L³ and L⁴ is a covalent bond, then u is the
 integer 1 and R⁵ is other than a hydrogen atom;
 and the salts, solvates, hydrates and N-oxides thereof.

20. (currently amended) A process according to Claim 1 for the preparation of
 compounds of formula (1d):



wherein

g is the integer 1, 2, 3 or 4;

~~R¹⁶, is an atom or group~~ -L³(Alk²)_tL⁴(R⁵)_u;

L³ and L⁴ are, independently, a covalent bond, an -O- or -S- atom, or a -C(O)-,
 -C(O)O-, -OC(O)-, -C(S)-, -S(O)-, -S(O)₂-, -N(R⁸)-, -CON(R⁸)-, -OC(O)N(R⁸)-, -CSN(R⁸),
 -N(R⁸)CO-, -N(R⁸)C(O)O-, -N(R⁸)CS-, -S(O)₂N(R⁸)-, -N(R⁸)S(O)₂-, -N(R⁸)O-, -ON(R⁸)-,
 -N(R⁸)N(R⁸)-, -N(R⁸)CON(R⁸)-, -N(R⁸)CSN(R⁸)-, or -N(R⁸)SO₂N(R⁸)- group;

R⁸ is a hydrogen atom or an optionally substituted straight or branched C₁-6alkyl
 group;

t is zero or the integer 1;

u is an integer 1, 2 or 3;

Alk² is an optionally substituted aliphatic chain or an optionally substituted heteroaliphatic chain containing one to four -O- or -S- atoms or -C(O)-, -C(O)O-, -OC(O)-, -C(S)-, -S(O)-, -S(O)₂-, -N(R⁸)-, -CON(R⁸)-, -OC(O)N(R⁸)-, -CSN(R⁸), -N(R⁸)CO-, -N(R⁸)C(O)O-, -N(R⁸)CS-, -S(O)₂N(R⁸)-, -N(R⁸)S(O)₂-, -N(R⁸)O-, -ON(R⁸)-, -N(R⁸)N(R⁸)-, -N(R⁸)CON(R⁸)-, -N(R⁸)CSN(R⁸)-, or -N(R⁸)SO₂N(R⁸)- groups that interrupt or are at the terminus of the aliphatic chain;

R⁵ is a hydrogen or halogen atom or an optionally substituted C₁-6alkyl, optionally substituted C₃-8cycloalkyl, -OR⁶, -SR⁶, -NR⁶R⁷, -NO₂, -CN, -CO₂R⁶, -SO₃H, -SOR⁶, -SO₂R⁶, -SO₃R⁶, -OCO₂R⁶, -CONR⁶R⁷, -OCONR⁶R⁷, -CSNR⁶R⁷, -COR⁶, -OCOR⁶, -N(R⁶)COR⁷, -N(R⁶)CSR⁷, -SO₂N(R⁶)(R⁷), -N(R⁶)SO₂R⁷, N(R⁶)CON(R⁷)(R¹⁹), -N(R⁶)CSN(R⁷)(R¹⁹), or -N(R⁶)SO₂N(R⁷)(R¹⁹) group; and

R⁶, R⁷, and R¹⁹ are, independently, a hydrogen atom or an optionally substituted C₁-6alkyl or C₃-8cycloalkyl group;

provided that when t is zero and each of L³ and L⁴ is a covalent bond, then u is the integer 1 and R⁵ is other than a hydrogen atom;

and the salts, solvates, hydrates and N-oxides thereof.

21. (previously presented) A process according to Claim 1 for the preparation of:

ethyl (2S)-2-[(2-bromo-3-oxospiro[3.5]non-1-en-1-yl)amino]-3-{4-[(3,5-dichloroisonicotinoyl)amino]phenyl}propanoate;

and the salts, solvates, hydrates and N-oxides thereof.

22. (previously presented) A process according to Claim 1 for the preparation of:

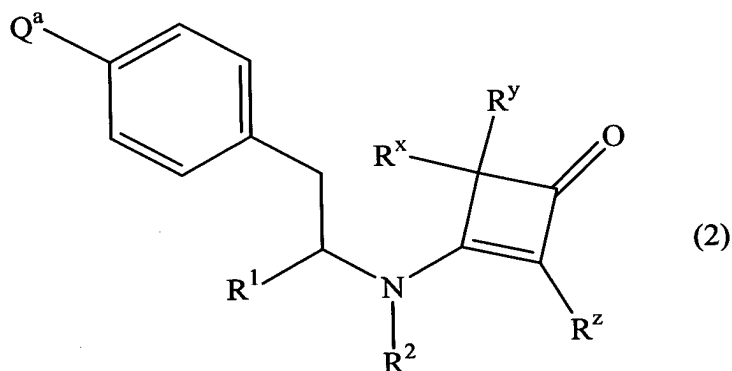
ethyl (2S)-2-(2-bromo-3-oxo-spiro[3.5]non-1-en-1-ylamino)-3-[4-([2,7]naphthyridin-1-ylamino)phenyl]propanoate;

and the salts, solvates, hydrates and N-oxides thereof.

23. (previously presented) A process according to Claim 1 for the preparation of:
 ethyl (2*S*)-2-[(2-isopropylsulfanyl-3-oxo-7-oxa-spiro[3.5]non-1-en-1-yl)amino]-3-[4-
 ([2,7]naphthyridin-1-ylamino)phenyl]propanoate;
 and the salts, solvates, hydrates and N-oxides thereof.

24. (previously presented) A process according to Claim 1 for the preparation of:
 2-hydroxyethyl (2*S*)-2-(2-bromo-3-oxo-spiro[3.5]non-1-en-1-ylamino)-3-{4-[(3,5-
 dichloroisonicotinoyl)amino]phenyl}propanoate;
 and the salts, solvates, hydrates and N-oxides thereof.

25. (currently amended) A compound of formula (2):



wherein:

R^1 is a carboxylic acid ($-\text{CO}_2\text{H}$) or ~~a derivative or biostere thereof~~ an acyclic or cyclic carboxylic acid ester, an amide, tetrazole, phosphonic acid, phosphinic acid, sulphonic acid, sulphinic acid, boronic acid, or an acylsulphonamide group;

R^2 is a hydrogen atom or a C_{1-6} alkyl group;

R^x , R^y and R^z , which may be the same or different, are each ~~an atom or group~~
 ~~$-\text{L}^1(\text{Alk}^1)_n(\text{R}^3)_v$ in which~~, or R^z is $-\text{L}^1(\text{Alk}^1)_n(\text{R}^3)_v$ and R^x and R^y are joined together to
form an optionally substituted spiro linked cycloaliphatic or heterocycloaliphatic group;

L^1 is a covalent bond or ~~a linker atom or group~~ an $-\text{O}-$, $-\text{S}-$, or $-\text{Se}-$ atom or an $-\text{C}(\text{O})-$, $-\text{C}(\text{O})\text{O}-$, $-\text{OC}(\text{O})-$, $-\text{C}(\text{S})-$, $-\text{S}(\text{O})-$, $-\text{S}(\text{O})_2-$, $-\text{N}(\text{R}^8)-$, $-\text{CON}(\text{R}^8)-$, $-\text{OC}(\text{O})\text{N}(\text{R}^8)-$, $-\text{CSN}(\text{R}^8)-$,

-N(R⁸)CO-, -N(R⁸)C(O)O-, -N(R⁸)CS-, -S(O)₂N(R⁸)-, -N(R⁸)S(O)₂-, -N(R⁸)O-, -ON(R⁸)-, -N(R⁸)N(R⁸)-, -N(R⁸)CON(R⁸)-, -N(R⁸)CSN(R⁸)-, or -N(R⁸)SO₂N(R⁸)- group;

R⁸ is a hydrogen atom or an optionally substituted straight or branched C₁₋₆alkyl group;

Alk¹ is an optionally substituted aliphatic chain or an optionally substituted heteroaliphatic chain containing one to four -O- or -S- atoms or -C(O)-, -C(O)O-, -OC(O)-, -C(S)-, -S(O)-, -S(O)₂-, -N(R⁸)-, -CON(R⁸)-, -OC(O)N(R⁸)-, -CSN(R⁸), -N(R⁸)CO-, -N(R⁸)C(O)O-, -N(R⁸)CS-, -S(O)₂N(R⁸)-, -N(R⁸)S(O)₂-, -N(R⁸)O-, -ON(R⁸)-, -N(R⁸)N(R⁸)-, -N(R⁸)CON(R⁸)-, -N(R⁸)CSN(R⁸)-, or -N(R⁸)SO₂N(R⁸)- groups that interrupt or are at the terminus of the aliphatic chain;

R³ is a hydrogen or halogen atom or group selected from -OR^{3a} [~~where R^{3a} is a hydrogen atom or an optionally substituted straight or branched C₁₋₆alkyl group or C₃₋₈cycloalkyl group~~], -SR^{3a}, -CN and an optionally substituted cycloaliphatic, heterocycloaliphatic, polycycloaliphatic, heteropolycycloaliphatic, aromatic or heteroaromatic group;

R^{3a} is a hydrogen atom or an optionally substituted straight or branched C₁₋₆alkyl group or C₃₋₈cycloalkyl group;

n is zero or the integer 1; and

v is the integer 1, 2 or 3;

provided that when n is zero and L¹ is a covalent bond, v is the integer 1;

~~or R^z is an atom or group as previously defined and R^x and R^y are joined together to form an optionally substituted spiro-linked cycloaliphatic or heterocycloaliphatic group;~~

Q^a is a group -N(R⁴)H;

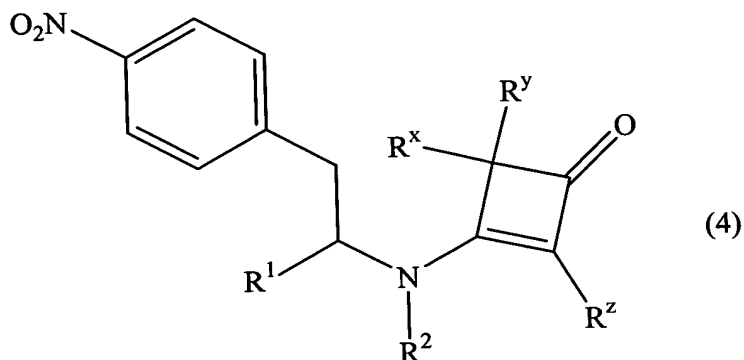
R⁴ is a hydrogen atom or an optionally substituted straight or branched C₁₋₆alkyl group;

and the salts, solvates, hydrates and N-oxides thereof.

26. (original) A compound according to Claim 25 which is:

3-(4-aminophenyl)-2(S)-(3-oxo-7-oxaspiro[3.5]non-1-en-1-ylamino)-propionic acid hydroxyethyl ester.

27. (currently amended) A compound of formula (4):



wherein:

R^1 is a carboxylic acid ($-\text{CO}_2\text{H}$) or ~~a derivative or biostere thereof~~ an acyclic or cyclic carboxylic acid ester, an amide, tetrazole, phosphonic acid, phosphinic acid, sulphonic acid, sulphinic acid, boronic acid, or an acylsulphonamide group;

R^2 is a hydrogen atom or a C_{1-6} alkyl group;

R^x , R^y and R^z , which may be the same or different, are each ~~an atom or group~~ $-\text{L}^1(\text{Alk}^1)_n(\text{R}^3)_v$ in which, or R^z is $-\text{L}^1(\text{Alk}^1)_n(\text{R}^3)_v$ and R^x and R^y are joined together to form an optionally substituted spiro linked cycloaliphatic or heterocycloaliphatic group;

L^1 is a covalent bond or ~~a linker atom or group~~ an $-\text{O}-$, $-\text{S}-$, or $-\text{Se}-$ atom or an $-\text{C}(\text{O})-$, $-\text{C}(\text{O})\text{O}-$, $-\text{OC}(\text{O})-$, $-\text{C}(\text{S})-$, $-\text{S}(\text{O})-$, $-\text{S}(\text{O})_2-$, $-\text{N}(\text{R}^8)-$, $-\text{CON}(\text{R}^8)-$, $-\text{OC}(\text{O})\text{N}(\text{R}^8)-$, $-\text{CSN}(\text{R}^8)-$, $-\text{N}(\text{R}^8)\text{CO}-$, $-\text{N}(\text{R}^8)\text{C}(\text{O})\text{O}-$, $-\text{N}(\text{R}^8)\text{CS}-$, $-\text{S}(\text{O})_2\text{N}(\text{R}^8)-$, $-\text{N}(\text{R}^8)\text{S}(\text{O})_2-$, $-\text{N}(\text{R}^8)\text{O}-$, $-\text{ON}(\text{R}^8)-$, $-\text{N}(\text{R}^8)\text{N}(\text{R}^8)-$, $-\text{N}(\text{R}^8)\text{CON}(\text{R}^8)-$, $-\text{N}(\text{R}^8)\text{CSN}(\text{R}^8)-$, or $-\text{N}(\text{R}^8)\text{SO}_2\text{N}(\text{R}^8)-$ group;

R^8 is a hydrogen atom or an optionally substituted straight or branched C_{1-6} alkyl group; ;

Alk^1 is an optionally substituted aliphatic chain or an optionally substituted heteroaliphatic chain containing one to four $-\text{O}-$ or $-\text{S}-$ atoms or $-\text{C}(\text{O})-$, $-\text{C}(\text{O})\text{O}-$, $-\text{OC}(\text{O})-$,

-C(S)-, -S(O)-, -S(O)₂-, -N(R⁸)-, -CON(R⁸)-, -OC(O)N(R⁸)-, -CSN(R⁸), -N(R⁸)CO-,
-N(R⁸)C(O)O-, -N(R⁸)CS-, -S(O)₂N(R⁸)-, -N(R⁸)S(O)₂-, -N(R⁸)O-, -ON(R⁸)-,
-N(R⁸)N(R⁸)-, -N(R⁸)CON(R⁸)-, -N(R⁸)CSN(R⁸)-, or -N(R⁸)SO₂N(R⁸)- groups that
interrupt or are at the terminus of the aliphatic chain;

R³ is a hydrogen or halogen atom or group selected from -OR^{3a} [~~where R^{3a} is a
hydrogen atom or an optionally substituted straight or branched C₁₋₆alkyl group or C₃₋₈
cycloalkyl group~~], -SR^{3a}, -CN and an optionally substituted cycloaliphatic,
heterocycloaliphatic, polycycloaliphatic, heteropolycycloaliphatic, aromatic or
heteroaromatic group; ;

R^{3a} is a hydrogen atom or an optionally substituted straight or branched C₁₋₆alkyl
group or C₃₋₈cycloalkyl group;

n is zero or the integer 1; and

v is the integer 1, 2 or 3;

provided that when n is zero and L¹ is a covalent bond, v is the integer 1;

~~or R^Z is an atom or group as previously defined and R^{*} and R^Y are joined together to form an
optionally substituted spiro-linked cycloaliphatic or heterocycloaliphatic group;~~
and the salts, solvates, hydrates and N-oxides thereof.

28. (original) A compound according to Claim 27 which is:

3-(4-nitrophenyl)-2(S)-(3-oxo-7-oxaspiro[3.5]non-1-en-1-ylamino)propionic acid
hydroxyethyl ester.